## **Claims**

We claim:

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A compound of the structure

q Y A E T  $R^4$ 

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wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>1</sup>, C(R<sup>2</sup>)(R<sup>3</sup>), NR<sup>5</sup>, CH, O and S;

q is an integer of from 3 to 10;

A is selected from the group consisting of O, S, C(R<sup>16</sup>)(R<sup>17</sup>) and NR<sup>6</sup>;

E is selected from the group consisting of CH<sub>2</sub>, O, S, and

 $NR^7$ ;

J is selected from the group consisting of O, S and NR<sup>8</sup>;

T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of from 0 to 3;

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M is selected from the group consisting of  $\overline{C}(R^9)(R^{10})$  and  $(CH_2)_u$ , wherein u is an integer of from 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

X is selected from the group consisting of CO<sub>2</sub>B, PO<sub>3</sub>H<sub>2</sub>,

SO<sub>3</sub>H, SO<sub>2</sub>NH<sub>2</sub>, SO<sub>2</sub>NHCOR<sup>12</sup>, OPO<sub>3</sub>H<sub>2</sub>, C(O)NHC(O)R<sup>13</sup>,

C(O)NHSO<sub>2</sub>R<sup>14</sup>, hydroxyl, tetrazolyl and hydrogen;

W is selected from the group consisting of C, CR<sup>15</sup> and N; and

B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup>, R<sup>14</sup>, R<sup>15</sup>, R<sup>16</sup> and R<sup>17</sup> at each occurrence are independently selected from the group consisting of

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A.A.

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hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C1-C3 alkyl)-C(O)(C1-C3 alkyl), -NHC(O)N(C1-C3 alkyl)C(O)NH(C1-C3alkyl), -NHC(O)NH(C1-C6alkyl), -NHSO2(C1-C3 alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>- $C_3$ )amino, $-C(O)O-(C_1-C_3)$ alkyl,  $-C(O)NH-(C_1-C_3)$ alkyl,  $-C(O)N(C_1-C_3)$ alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein B,  $R^1$ ,  $R^2$ ,  $R^3$ ,  $R^4$ ,  $R^5$ ,  $R^6$ ,  $R^7$ ,  $R^8$ ,  $R^9$ ,  $R^{10}$ ,  $R^{11}$ ,  $R^{12}$ ,  $R^{13}$ , R<sup>14</sup>, R<sup>15</sup>, R<sup>16</sup> and R<sup>17</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; and wherein when M is C(R9)(R10), R9 and R10 taken together may form a ring;

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taken together may form a ring;
or a pharmaceutically acceptable salt thereof;

and wherein when A is NR<sup>6</sup> and at least one Y is CR<sup>1</sup>, R<sup>1</sup> and R<sup>6</sup>

or a pharmaceutically acceptable salt thereof; with the proviso that when A is  $C(R^{16})(R^{1})$ , E is not  $NR^{7}$ .

25 2. A compound of claim 1 wherein

A is NR<sup>6</sup>;
E is NR<sup>7</sup>;
J is O;
M is C(R<sup>9</sup>)(R<sup>10</sup>);

q is 4 or 5; T is (CH<sub>2</sub>)<sub>b</sub> wherein b is 0; X is CO<sub>2</sub>B;

W is C or CR<sup>15</sup>;

R<sup>4</sup> is selected from the group consisting of aryl, alkylaryl, aralkyl, heterocyclyl, alkylheterocyclyl and heterocyclylalkyl; and R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup> and R<sup>15</sup> are independently selected from the group consisting of hydrogen and lower alkyl.

3. A compound of claim 1 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

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A compound of the structure

$$R^9$$
 $R^9$ 
 $R^9$ 

1.5

wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>1</sup>, C(R<sup>2</sup>)(R<sup>3</sup>), NR<sup>5</sup>, CH, O and S;

q is an integer of from 3 to 7;

T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

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W is selected from the group consisting of C, CR<sup>15</sup> and N; and

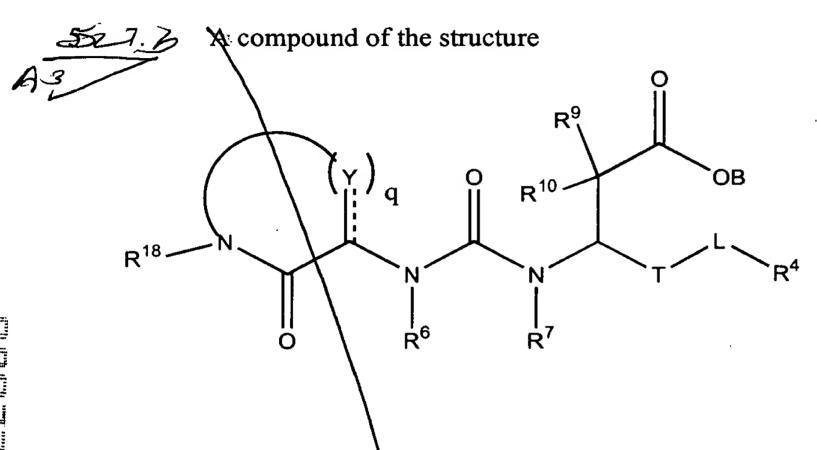
B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>15</sup> are independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

Heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N( $C_1$ - $C_3$  alkyl)C(O)NH( $C_1$ - $C_3$ alkyl), -NHC(O)NH( $C_1$ - $C_6$  alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-C3)amino, -C(O)O-(C1-C3)alkyl, -C( $\Diamond$ )NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, 5 -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SQ<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and 10 -C(O)NH(benzyl) groups; wherein B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>15</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; 15 and wherein R and R and R taken together may form a ring; and wherein when at least one Y is CR<sup>1</sup>, R<sup>1</sup> and R<sup>6</sup> taken together may form a ring; or a pharmaceutically acceptable salt thereof. 20 A compound of claim 4 wherein q is 4 or 5; W is C or CR<sup>15</sup>; T is  $(CH_2)_b$  wherein b is 0; 25 L is  $(CH_2)_n$  wherein n is 0; R<sup>4</sup> is selected from the group consisting of aryl, alkylaryl, aralkyl, heterocyclyl, alkylheterocyclyl and heterocyclylalkyl; and

R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup> and R<sup>15</sup> are independently selected from the

group consisting of hydrogen and lower alkyl.

6. A compound of claim 4 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.



wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>1</sup>, C(R<sup>2</sup>)(R<sup>3</sup>), NR<sup>5</sup>, CH, O and S;

q is an integer of from 2 to 5;

T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and

B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>9</sup> and R<sup>10</sup> are independently selected from the group consisting of hydrogen, halogen, halkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

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heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl),

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-NHC(O)N( $C_1$ - $C_3$  alkyl)C(O)NH( $C_1$ - $C_3$ alkyl), -NHC(O)NH( $C_1$ - $C_6$  alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-C3)amino, -C(O)O-(C1-C3)alkyl,  $-C(Q)NH-(C_1-C_3)alkyl, -C(Q)N(C_1-C_3)alkyl)_2, -CH=NOH, -PO_3H_2,$ **5** . -OPOH2, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy,\arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and 10 -C(O)NH(benzyl) groups; wherein B,  $\mathbb{R}^{1}$ ,  $\mathbb{R}^{2}$ ,  $\mathbb{R}^{3}$ ,  $\mathbb{R}^{4}$ ,  $\mathbb{R}^{5}$ ,  $\mathbb{R}^{6}$ ,  $\mathbb{R}^{7}$ ,  $\mathbb{R}^{9}$ ,  $\mathbb{R}^{10}$ ,  $\mathbb{R}^{11}$  and  $\mathbb{R}^{18}$  are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; 15 and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring; and wherein when at least one Y is CR1, R1 and R6 taken together may form a ring; or a pharmaceutically acceptable salt thereof. 20

8. A compound of claim 7 wherein R<sup>18</sup> is selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, cycloalkyl, alkylheterocyclyl, heterocyclylalkyl and heterocyclyl;

T is (CH<sub>2</sub>)<sub>b</sub> wherein b is 0;

L is  $(CH_2)_n$  wherein n is 0; Y is selected from the group consisting of  $CR^1$  and  $C(R^2)(R^3)$  and q is 2 or 3.

9. A compound of claim 7 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

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10. A compound of claim 7 wherein

is selected from the group consisting of

$$\mathbb{R}^{22}$$

$$\mathbb{R}^{18}$$
and

wherein R<sup>19</sup>, R<sup>20</sup>, R<sup>21</sup> and R<sup>28</sup> at each occurrence are independently selected from the group consisting of halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -OH, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, atolyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, atalkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl,

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sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups;

R<sup>18</sup> is selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

R<sup>22</sup> is selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups;

c is an integer of zero to two;

d is an integer of zero to three;
e is an integer of zero to four; and
i is an integer of zero to two.

11. The compound of claim 7 wherein R<sup>18</sup> is aralkyl;

R<sup>4</sup> is aryl;

T is  $(CH_2)_b$  where b is zero;

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L is (CH<sub>2</sub>)<sub>n</sub> where n is zero; and,

B, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup> and R<sup>10</sup> are each independently hydrogen.

a compound of the structure

R<sup>10</sup>  $k^6$ 

> wherein T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of from 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

g is an integer of from 0 to 7; and

B,  $R^4$ ,  $R^9$ ,  $R^{10}$  and  $R^{23}$  at each occurrence are independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C( $\bigcirc$ )N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl\heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-

C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and

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-C(O)NH(benzyl) groups;

R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkerlyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterodycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl,\aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

wherein B, R<sup>4</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>18</sup> and R<sup>23</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; and wherein R and R and R taken together may form a ring;

or a pharmaceutically acceptable salt thereof.

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13. A compound of claim 12 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

14. A compound of the structure

$$R^{24}$$
 $R^{18}$ 
 $R$ 

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wherein h is an integer of zero to five;

B, R<sup>9</sup>, R<sup>10</sup>, R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl,

-CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

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heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N( $C_1$ - $C_3$  alkyl)C(O)NH( $C_1$ - $C_3$ alkyl), -NHC(O)NH( $C_1$ - $C_6$  alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>- $C_3$ )alkyl,  $-C(O)N(C_1-C_3 \text{ alkyl})_2$ , -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; R<sup>27</sup>, at each occurrence, is independently selected from the group consisting of halogen, hydroxyl, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl),  $-NHC(O)N(C_1-C_3\ alkyl)C(O)NH(C_1-C_3\ alkyl),\ -NHC(O)NH(C_1-C_6\ alkyl),$ -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), -N(C<sub>1</sub>-C<sub>3</sub>alkyl)SO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub>alkyl), -N(C<sub>1</sub>-C<sub>3</sub>alkyl)SO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>- $C_3$ )amino,  $-C(O)O-(C_1-C_3)$ alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; R<sup>6</sup>, R<sup>7</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, · alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy,

carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl,

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cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and, R<sup>26</sup> is selected from the group consisting of hydrogen, alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, alkoxycarbonyl, heterocycloyl, carboxy,  $-C(O)O-(C_1-C_3)$  alkyl,  $-C(O)NH-(C_1-C_3)$  alkyl,  $-C(O)N(C_1-C_3)$  alkyl)<sub>2</sub>,  $-C(O)N(C_1-C_3)$ PO<sub>3</sub>H<sub>2</sub>, haloalkyl, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, biaryl, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO2-(C1-C3 alkyl), sulfonamido, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein B,  $R^6$ ,  $R^7$ ,  $R^9$ ,  $R^{10}$ ,  $R^{18}$ ,  $R^{24}$ ,  $R^{25}$ ,  $R^{26}$  and  $R^{27}$  are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein R\square and R<sup>24</sup> taken together may form a ring; R<sup>24</sup> and R<sup>25</sup>\taken together may form a ring; R<sup>25</sup> and R<sup>26</sup> taken together may form a ring; and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring; or a pharmaceutically acceptable salt thereof.

- 15. The compound of claim 14 wherein B, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>24</sup>, R<sup>25</sup> and R<sup>26</sup> are each independently hydrogen and R<sup>18</sup> is substituted or unsubstituted aralkyl.
- 16. A compound of claim 14 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

17.\A compound of the structure

$$(Z)_z$$
 $(R^{29})_k$ 
 $R^{10}$ 
 $(R^{29})_k$ 
 $(R^{29})_k$ 

wherein Z, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>30</sup>, C(R<sup>31</sup>)(R<sup>32</sup>), NR<sup>33</sup>, CH, O and S;

z is an integer of from 3 to 6;

k is an integer of from 0 to 5;

T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of from 0 to 3;

L is selected from the group consisting of O,  $NR^{11}$ , S, and  $(CH_2)_n$  wherein n is an integer of 0 or 1;

R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup>, R<sup>18</sup> and R<sup>33</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

B, R<sup>4</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>30</sup>, R<sup>31</sup> and R<sup>32</sup> at each occurrence are independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -OH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl),

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-NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-C3)amino, -C(O)O-(C1-C3)alkyl, -C(O)NH-(C1-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl\ diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO2-(C1-C3 alkyl), -SO3-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; and

R<sup>29</sup>, at each occurrence, is independently selected from the group consisting of halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)- $C(O)(C_1-C_3 \text{ alk})$ , -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-C3)amino, -C(O)O-(C1- $C_3$ )alkyl,  $-C(O)NH^1(C_1-C_3)$ alkyl,  $-C(O)N(C_1-C_3)$ alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, białyl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups;

wherein B, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>18</sup>, R<sup>29</sup>, R<sup>30</sup>, R<sup>31</sup>, R<sup>32</sup> and R<sup>33</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group;

wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring;

or a pharmaceutically acceptable salt thereof.

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- 18. A compound of claim 17 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.
- 19. The compound of claim 17 wherein z is three or four.

## 20. A compound of the structure

hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy,

thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF3, -SH, -OH,

-CO<sub>2</sub>H, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl,

wherein R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of

carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub>

 $alkyl)C(O)NH(C_1-C_3alkyl), \ -NHC(O)NH(C_1-C_6alkyl), \ -NHSO_2(C_1-C_3alkyl), \ -NHSO_2(C_1-$ 

alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-

C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub>

alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy,

carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl,

cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl,

diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl,

heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl),

sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; and

R<sup>18</sup> and R<sup>34</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde,

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carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and –C(O)NH(benzyl) groups;

wherein R<sup>18</sup>, R<sup>24</sup>, R<sup>25</sup> and R<sup>34</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group;

and wherein  $R^{24}$  and  $R^{25}$  taken together may form a ring; with the proviso that when  $R^{24}$  and  $R^{25}$  taken together form a ring, the ring formed is not benzene.

21. A compound of claim 20 wherein R<sup>34</sup> is hydrogen;

R<sup>18</sup> is aralkyl; and R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of hydrogen, lower alkyl, and lower alkyl wherein R<sup>24</sup> and R<sup>25</sup> taken together may form a ring.

22. A compound of claim 20 of the structure

$$\mathbb{R}^{24}$$

$$\mathbb{R}^{25}$$

$$\mathbb{R}^{25}$$

$$\mathbb{R}^{25}$$

$$\mathbb{R}^{34}$$

wherein R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -SH, -OH, -CO<sub>2</sub>H, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub>

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alkyl), -NHSO2(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)2, -CH=NOH, -PO3H2, -OPO3H2, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; R<sup>34</sup> is selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and, R<sup>35</sup>, at each occurrence, is independently selected from the group consisting of halogen, hydroxyl, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl,  $-\tilde{SO}_2$ - $(\tilde{C}_1$ - $\tilde{C}_3$  alkyl), -SO<sub>3</sub>- $(C_1$ - $C_3$  alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein R<sup>24</sup>, R<sup>25</sup>, R<sup>34</sup> and R<sup>35</sup> are unsubstituted or substituted with

at least one electron donating or electron withdrawing group; and,

m is an integer of from 0 to 5.

- 5 23. A compound of claim 22 wherein R<sup>34</sup> is hydrogen;
  - m is an integer of one to three and R<sup>35</sup> at each occurrence is selected from the group consisting of alkyl, halogen, alkoxy, haloalkyl, sulfonyl, -OH and -CN.
  - 24. A compound of claim 20 selected from the group consisting of
- 5-(2-chlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-fluorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-fluorobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-benzyl-6-methyl-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-benzyl-3,5-dihydro[1,3]oxazolo[4,5-
- c]pyridine-2,4-dione, 5-(2,5-dimethylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4
  - dione, 5-(2-methylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-
  - dichlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-methoxybenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,5-difluorobenzyl)-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(methylthio)benzyl]-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-fluorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-methoxybenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3,5-bis(trifluoromethyl)benzyl]-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-tert-butylbenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-chlorobenzyl)-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(trifluoromethyl)benzyl]-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-bromobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,4-dichlorobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-methylbenzyl)-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methoxybenzyl)-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[4-(trifluoromethyl)benzyl]-3,5-

dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-methylbenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(pyridin-2-ylmethyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-difluorobenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-difluorobenzyl)-3,5-5 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(trifluoromethoxy)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[4-(trifluoromethoxy)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(trifluoromethyl)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-methoxybenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,3-dichlorobenzyl)-3,5-10 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,5-dimethylbenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-pentyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-dichlorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-ethyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-butyl-5-(2-chlorobenzyl)-3,5-15 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(trifluoromethyl)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dichlorobenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-fluorobenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methylbenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-7-methyl-3,5-20 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-5,6,7,8-tetrahydro-2Hcyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 7-methyl-5-[4-(methylsulfonyl)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4methoxybenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7propyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 4-[(2,4-dioxo-2,3-25 dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]-N,N-dimethylbenzenesulfonamide, 5-(mesitylmethyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-3,5,6,7,8,9-hexahydro[1,3]oxazolo[4,5-c]quinoline-2,4-dione, 5-(2-chlorobenzyl)-7-ethyl-6methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(methylthio)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 2-[(2,4-dioxo-2,3-dihydro[1,3]oxazolo[4,5-30

c]pyridin-5(4H)-yl)methyl]-N,N-dimethylbenzenesulfonamide, 5-(2,6-dimethoxybenzyl)-3,5-

dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(trifluoromethoxy)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-6,7-dimethyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(methylsulfonyl)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chloro-2-methoxybenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-5,6,7,8,9,10-hexahydro-5 2H-cyclohepta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-[2-(difluoromethoxy)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-[(1R)-1-phenylethyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-7-propyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(methylsulfonyl)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dimethylbenzyl)-3,5-10 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 3-chloro-2-[(2,4-dioxo-2,3dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]benzonitrile, 5-(2-chloro-6methylbenzyl)-6,7-dimethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 2-[(2,4-dioxo-2,3-dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]benzonitrile, 5-(2-chloro-6methoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-15 (methylthio)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7cyclopropyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-chlorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dichlorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-(4-methylbenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,5-dimethoxybenzyl)-7-methyl-3,5-20 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-difluorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(methylsulfonyl)benzyl]-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-fluoro-6-methoxybenzył)-7-methyl-3,5-25 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methoxybenzyl)-7-propyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-chloro-2-fluorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-isopropyl-3,5dinydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-fluoro-2-methylbenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-[(1S)-1-phenylethyl]-3,5-30 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-isopropoxybenzyl)-7-methyl-

3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-acetyl-2-methoxybenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-[2-fluoro-6-(trifluoromethyl)benzyl]-7methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methylbenzyl)-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-(2-chloro-5 6-ethoxybenzyl)-7-ethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6propoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6isobutoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)dione, 5-(2-chloro-6-isopropoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-10 2,4-dione, 5-[2-chloro-6-(2,2,2-trifluoroethoxy)benzyl]-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-[2-chloro-6-(2-methoxyethoxy)benzyl]-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-(2-chloro-6-ethoxybenzyl)-6,7-dimethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-15 6-ethoxybenzyl)-7-ethyl-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2chlorobenzyl)-7-ethyl-3,5-dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-7-propyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-7-cyclopropyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-propoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-20 methoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5ethoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(piperidin-1-ylsulfonyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(pyrrolidin-1-ylsulfonyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-25 c]pyridine-2,4-dione, 5-[2-chloro-6-(cyclopentylmethoxy)benzyl]-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(benzyloxy)-6-chlorobenzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,3-dichloro-6-ethoxybenzyl)-5,6,7,8tetranydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-[2-chloro-5-(trifluoromethyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione and 5-30 (2-chloro-5-fluorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione.

- 25. A compound selected from the group consisting of (3S)-3-[({[2-methyl-4-(2-methylpropyl)-6-oxo-1-(phenylmethyl)-1,6-dihydro-5-pyrimidinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenylphen
- benzodioxol-5-yl)-3-[({[2-oxo-1-(phenylmethyl)-4-propyl-1,2-dihydro-3-pyridinyl]amino}carbonyl)amino]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-ethyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2-oxo-4-propyl-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid,
- 10 (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid,

  (3S)-3-{[({6-methyl-2-oxo-1-(phenylmethyl)-4-[(phenylmethyl)oxy]-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2,4-dimethyl-6-oxo-1,6-dihydro-5-yrimidinyl}amino)carbonyl]amino}-
- 3-(4-methylphenyl)propanoic acid, (3S)-3-{[({4-amino-1-[(2-chlorophenyl)methyl]-6-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-
- (3,4-dimethylphenyl)propanoic acid, (3S)-3-{[({4-amino-1-[(2-chlorophenyl)methyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-4-(1,4-oxazinan-4-yl)-2-oxo-1,2-dihydro-3-
- pyridinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-2-oxo-4-(propylamino)-1,2-dihydro-3-pyridinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-bromophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl
- dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-methyl-4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2-oxo-4-phenyl-1,2-dihydro-3-

- pyrazinyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-
- pyridinyl}amino)carbonyl]amino}-3-(3,5-dimethylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(3-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-(methyloxy)phenyl]propanoic acid, (3S)-3-[3,5-bis(methyloxy)phenyl]-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}
- chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-quinolinyl}amino)carbonyl]amino}-3(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-(trifluoromethyl)phenyl]propanoic
  acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4({ethyl[(ethylamino)carbonyl]amino}carbonyl)amino]-2-oxo-1,2-dihydro-3-
- pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({4-(1-azetanyl)-1-[(2-chlorophenyl)methyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-4-({2-[(2-{[2-(methyloxy)ethyl]oxy}ethyl)oxy]ethyl}oxy)-2-oxo-1,2-dihydro-3-pyridinyi]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-
- 30 {[({1-[(2-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-fluorophenyl)methylphenyl)propanoic acid, (3S)-3-({1-[(2-fluorophenyl)methylphenyl)propanoic acid, (3S)-3-({1-[(2-fluorophenyl)methylphenyl

chloro-6-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2chlorophenyl)methyl]-5-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-(trifluoromethyl)phenyl)methyl)-1,2 dihydro-3-pyridinyl)amino)carbonyl)amino)propanoic 5 acid, (3S)-3-((((1-((2-chlorophenyl)methyl)-2-oxo-1,2-dihydro-3pyridinyl)amino)carbonyl)amino)-3-(4-methylphenyl)propanoic acid, (3S)-3-((((1-((2fluorophenyl)methyl)-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4methylphenyl)propanoic acid, (3S)-3-((((1-((2-bromophenyl)methyl)-2-oxo-1,2-dihydro-3pyridinyl)amino)carbonyl)amino)-3-(4-methylphenyl)propanoic acid, (3S)-3-((((1-((2,4-10 dichlorophenyl)methyl)-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4methylphenyl)propanoic acid, (3S)-3-((((1-((2-chloro-6-fluorophenyl)methyl)-2-oxo-1,2dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4-methylphenyl)propanoic acid, (3S)-3-((((1-((2-chlorophenyl)methyl)-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4-trifluoromethyl)oxy)phenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-2-15 oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, 4- $\{[3-[(\{[(1S)-2-carboxy-1-(4-methylphenyl)ethyl]amino\} carbonyl)amino]-1-(2-chlorobenzyl)-1-(4-methylphenyl)ethyl]amino\} carbonyl)amino]-1-(4-methylphenyl)ethyl]amino\} carbonyl)amino]-1-(4-methylphenyl)ethyl]amino\} carbonyl)amino]-1-(4-methylphenyl)ethyl]amino} carbonyl)amino]-1-(4-methylphenyl)ethyllamino]-1-(4-methylphenyl)eth$ 2-oxo-1,2-dihydropyridin-4-yl]amino}benzoic acid, (3S)-3-{[({1-(2-chlorobenzyl)-4-[(2,2dimethylpropanoyl)amino]-2-oxo-1,2-dihydropyridin-3-yl}amino)carbonyl]amino}-3-(4methylphenyl)propanoic acid, (3S)-3-[({[4-{[(tert-butylamino)carbonyl]amino}-1-(2-20 chlorobenzyl)-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4methylphenyl)propanoic acid, (3S)-3-[({[1-(2-cyanobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2,3dihydro-1,4-benzodioxin-6-yl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-25 oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(7-methoxy-1,3-benzodioxol-5yl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(3-ethoxy-4-methoxyphenyl)propanoic acid, (3S)-3-[({[1-(2chlorobenzy!) 4 hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4dimethoxyphenyl)propanoic acid, (3S)-3-[({[1-(4-chlorobenzyl)-4-hydroxy-2-oxo-1,2-30 dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-

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(2-chloro-6-methoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
     yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
     4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-
     methylphenyl)propanoic acid, (3S)-3-[({[1-(2,6-difluorobenzyl)-4-hydroxy-2-oxo-1,2-
     dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-
 5
     (2-chloro-6-methoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
     yl]amino}carbonyl)amino]-3-(3,5-dimethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-
     chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-
     diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-
     dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-
10
     (2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-
     methoxy-4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-
      1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,5-dimethoxy-4-
     methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-
     dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-dimethylphenyl)propanoic acid, (3S)-3-
15
     [({[1-(2-chlorobenzyl)-5-ethyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
     yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[2-chloro-5-
     (trifluoromethyl)benzyl]-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl}amino)carbonyl]amino}-
      3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-4-hydroxy-2-
     oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-methylphenyl)propanoic acid,
20
     (3S)-3-[({[1-(2-chloro-6-methylbenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-
     yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
     4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-
     (4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2,6-dimethoxybenzyl)-4-hydroxy-2-oxo-1,2-
     dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-
25
     (2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-
     propoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-5-propyl-1,2-
     dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-
     (2 chlorobonzyl)-4-liydroxy-5,ó-áimethyl-2-oxo-1,2-dihydropyridin-3-
     yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
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4-hydroxy-2-oxo-5-propyl-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-

diethoxyphenyl)propanoic acid, (3S)-3-(3-butoxyphenyl)-3-[({[1-(2-chlorobenzyl)-4hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]propanoic acid, (3S)-3-{[({1-[2-chloro-5-(methylsulfonyl)benzyl]-4-hydroxy-2-oxo-1,2-dihydropyridin-3yl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(2-5 methoxyethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-dipropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(difluoromethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-10 diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methylbenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(3,4-diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6cyanobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-15 methylphenyl)propanoic acid, 3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2-naphthyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4diethoxyphenyl)propanoic acid (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-4-hydroxy-5methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-20 diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(4-methoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6methylbenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-25 yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-CKC-1,2-diliydropyridin-3-yi]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-30 cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-

- [({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2,3-dihydro-1-benzofuran-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-
- hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3,5-diethoxyphenyl)propanoic acid, (3S)-3-[({[5-chloro-1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-dihydropyridin-3-yl]amino}carbonyl
- [({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-propoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-phenylpropanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-
- yl]amino}carbonyl)amino]-3-(1,3-diethyl-2-oxo-2,3-dihydro-1H-benzimidazol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(trifluoromethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dihydroy-1-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dihydroy-1-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dihydroy-1-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)amino]-3-(3-isopropoxyphenyl)amino]-3-(3-isopropoxyphenyl)amino]-3-(3-isopropoxyphenyl)amino]-3-(3-isopropoxyp
- chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-5-cyclopropyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-5-cyclopropyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
- methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-6-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-

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tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-[3(cyclopropyloxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-[3(cyclopropylmethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3[3-(cyclopropylmethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3,5-dimethylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3[(difluoromethyl)oxy]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-

hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3 {3-[(1,1,2,2-tetrafluoroethyl)oxy]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-(1-ethyl-1H-indol-5-yl)propanoic acid and (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-[3-(diethylamino)phenyl]propanoic acid and pharmaceutical acceptable salts thereof.

26. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid and pharmaceutically acceptable salts thereof.

27. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid and pharmaceutically acceptable salts thereof.

28. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(diethylamino)phenyl]propanoic acid and pharmaceutically acceptable salts thereof.

- 29. A compound selected from the group consisting of (3S)-3-[({[1-(2-chloro-6ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic 5 acid; (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1Hcyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(6-methoxy-2-naphthyl)propanoic acid; (3S)-3-[({[1-(2chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-10 yl]amino}carbonyl)amino]-3-(3-methylphenyl)propanoic acid; (3S)-3-{[({1-[(2-chloro-6methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3yl}amino)carbonyl]amino}-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-{[({1-[(2chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3yl}amino)carbonyl]amino}-3-{3-[(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-15 {[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1Hcyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[methyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-20
- cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3
  [ethyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[ethyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-(1H-indol-5-yl)propanoic acid and pharmaceutically acceptable salts thereof.

methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-

3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-

yl}amino)carbonyl]amino}-3-{3-[methyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-

- 30. A pharmaceutical composition comprising:
  a compound of claim 1
  in a pharmaceutically acceptable carrier.
- 5 31. A method for selectively inhibiting  $\alpha_4\beta_1$  integrin binding in a mammal comprising administering to said mammal a therapeutic amount of a compound of claim 1.